

On the Quantum Circuit Complexity Equivalence

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Abstract

Nielsen [3] recently asked the following question: "What is the minimal size quantum circuit required to exactly implement a specified n -qubit unitary operation U , without the use of ancilla qubits?" Nielsen was able to prove that a lower bound on the minimal size circuit is provided by the length of the geodesic between the identity I and U , where the length is defined by a suitable Finsler metric on $SU(2^n)$. We prove that the minimum circuit size that simulates U is in linear relation with the geodesic length and simulation parameters, for the given Finsler structure F . As a corollary we prove the highest lower bound of $O(\frac{n^4}{p} d_{F_p}^2(I, U) L_{F_p}(I, \tilde{U}))$ and the lowest upper bound of $\Omega(n^4 d_{F_p}^3(I, U))$, for the standard simulation technique. Therefore, our results show that by standard simulation one can not expect a better than n^2 times improvement in the upper bound over the result from Nielsen, Dowling, Gu and Doherty [4]. Moreover, our equivalence result can be applied to the arbitrary path on the manifold including the one that is generated adiabatically.

1 Introduction

Quantum computation is inherently a process of continuous evolution of quantum states that has the potential to fundamentally change the notion of feasibly tractable computation. Only recently did researchers start to think how notions from the differential geometry [13] can be used to represent this process. Instead, quantum circuits, as an inherently discrete notion of computation, have been most commonly used to represent this continuous process. Any quantum operation, a unitary matrix $U \in SU(2^n)$, is an element of a Lie group, and a point on the $\mathcal{U} \equiv SU(2^n)$ manifold, whose tangent bundle can be endowed with the Finsler structure F , that effectively provides a measure of length for any path on the Finsler manifold (\mathcal{U}, F) . In particular, the paths that we are interested in are geodesics. These are locally and, under certain conditions, globally minimal length paths between any two points on the manifold. They are of particular interest, because if closely simulated they can give the smallest circuit complexity for the given unitary U .

The aim in the approach that we take here is to tackle the question about the complexity of the circuit necessary for the simulation of an arbitrary unitary gate. As the length of the geodesic for the particular unitary is its intrinsic property, ideally one would succeed in finding the minimum number of circuits necessary to implement the unitary by simulating exactly its geodesic. Therefore, the hope here is to learn about the circuit lower bounds by basically transforming the hard combinatorial optimization problems over large sets to the problems in continuous domain that can be solved with tools of differential geometry and the calculus of variations.

One of the first results that had the flavor of this transform was introduced by Mochon [5], who proved that in the discrete model and the analogous continuous model, i.e. the Hamiltonian oracle model, Oracle interrogation, the problem of computing XOR and Grover search have the same complexity. Moreover, Nielsen [3] and subsequently Nielsen, Dowling, Gu and Doherty [4] proved that for particularly chosen metric there is a polynomial equivalence between the geodesic length and number of gates necessary for the simulation. The lower bound for the minimum number of gates necessary for the simulation has been determined for exact simulation and the upper bound has been determined for the arbitrary precision. The metric chosen in [4] penalizes all those directions on the manifold that are not easily simulated by local gates, so that coefficients for stabilizer elements of Hamming weight greater than two bear high cost, i.e. have longer paths.

In this paper we prove the stronger result and show the exact upper and lower bound that determine the equivalence between the minimal number of gates in the standard circuit simulation and the length of the geodesic. Both upper and lower bound are determined by the simulation parameters and, of course, the length of the geodesic.

We consider the two cases. First: the simulation of the geodesic with set of gates \mathcal{G} that is exactly universal, and the case with approximately universal set of gates. With the exactly universal set of gates for any point $x_0 \in \mathcal{U}$ there is a gate in a set \mathcal{G} by which we can simulate exactly any point $x_1 \in \mathcal{U}$ in the ball of radius r centered around x_0 ; we denote this ball as $B_{x_0}^+(r)$. Under that assumption, we prove that the number of gates in the simulation of a geodesic may be upper and lower bounded by a linear factor in the length of geodesic and simulation parameters.

When the set of gates, \mathcal{G}_ϵ , is approximately universal, a single gate from this set can simulate the points in $B_{x_0}^+(r)$ only with some finite precision ϵ , and that will necessarily mean that the circuit that simulates the geodesic is doing so along the path that is not shorter than the actual geodesic, for that very point.

Our aim here is not to elaborate on the algorithm for the geodesic simulation but rather to prove the bounds that optimal simulation can achieve. We say optimal, because the set of gates \mathcal{G} that we first consider is much more powerful than any local and universal set of gates. Therefore the result that we present is the optimal result about complexity equivalence between discrete and continuous notions of computation. In particular, for the standard simulation model described in [4], we derive the highest lower bound and the lowest upper bound in the minimal circuit complexity that one can hope to achieve with the simulation of a geodesic.

2 Preliminaries

A quantum operation $U \in \mathcal{U}$ is a point on the manifold $\mathcal{U} \equiv SU(2^n)$ at some distance from identity $I \in \mathcal{U}$. The distance considered is the integral distance that is determined by the structure used on the manifold. In general that structure may be more general than Riemannian, i.e. it is called the Finsler structure $F(x, y)$. The restriction of a Finsler structure F to any specific tangent space $T_x\mathcal{U}$ with the origin at the point $x \in \mathcal{U}$ is called Minkowski norm on $T_x\mathcal{U}$. The second argument of the structure $F(x, y)$ is the velocity and its definition follows. Therefore a Finsler structure is basically a family of the smoothly varying Minkowski norms, one for each tangent space.

The defining properties of a non-negative real-valued structure $F(x, y)$ on \mathbb{R}^{4^n-1} are as follows:

- (1) it is C^∞ anywhere on \mathbb{R}^{4^n-1} except at $y = 0$;
- (2) it is positive homogeneous, i.e. $F(x, \lambda y) = \lambda F(x, y)$ for $\lambda > 0$;
- (3) the $(4^n - 1) \times (4^n - 1)$ matrix $\frac{\partial}{\partial y_i} \frac{\partial}{\partial y_j} [\frac{1}{2} F^2]$ is positive definite unless $y = 0$. As a consequence, one can derive positivity and triangular equality of Minkowski norms [6]. The structure $F(x, y)$ is usually denoted simply as $F(y)$.

For any $a, b \in \mathbb{R}_+$ we say that a map $\sigma : [a, b] \rightarrow \mathcal{U}$ is a piecewise C^∞ curve with velocity $y \equiv \frac{d\sigma}{dt} = \sum_i \frac{d\sigma_i}{dt} \frac{\partial}{\partial x_i} \in T_{\sigma(t)} \mathcal{U}$. The integral length of the curve σ , $L(\sigma)$, is defined as:

$$L_F(\sigma) = \int_a^b F(\sigma, \frac{d\sigma}{dt}) dt . \quad (1)$$

Since we are usually interested in minimum length curves for $x_0, x_1 \in \mathcal{U}$, we denote by $\Gamma(x_0, x_1)$ the collection of all piecewise C^∞ curves $\sigma : [a, b] \rightarrow \mathcal{U}$ such that $\sigma(a) = x_0$ and $\sigma(b) = x_1$. Similarly, the integral distance is defined as a map $d_F : \mathcal{U} \times \mathcal{U} \rightarrow [0, \infty)$:

$$d_F(x_0, x_1) = \inf_{\Gamma(x_0, x_1)} L_F(\sigma) \quad (2)$$

Using these definitions, one can show that the Finsler manifold (\mathcal{U}, d_F) satisfies the two axioms of a metric space: (1) positivity: $d_F(x_0, x_1) \geq 0$, where equality holds if and only if $x_0 = x_1$ and (2) the triangular inequality: $d_F(x_0, x_2) \leq d_F(x_0, x_1) + d_F(x_1, x_2)$. In general, the symmetric property of a distance does not need to hold, and therefore $d_F(x_0, x_1) \neq d_F(x_1, x_0)$.

3 Distortion Lemma

To establish the equivalence result, we introduce in this section the main tool of our analysis. The intuitive idea on which we build our results relies on the relation between the distances on the manifold and the distances on the tangent space of the manifold. While the former are introduced by the unitary gates and their complexity, the latter are defined by the appropriately defined distances between the Hamiltonians of gates used in the simulation. This will be proven useful in the sections below.

The lemma that follows is a slightly stronger result of a well-known and very useful fact from the differential geometry. Again, it relates distances on the manifold with the minimum and maximum distortion of the Euclidian norm on the tangent space over the compact set. Interested reader are encouraged to consult [6], an excellent and very elaborate reference on this subject.

Lemma 1 (Distortion Lemma). *Let (\mathcal{U}, F) be a Finsler manifold, and for any point $x \in \mathcal{U}$ let $\varphi : P_x \rightarrow \mathbb{R}^{4^n-1}$ be the local coordinate system diffeomorphism of a compact set P_x onto an open ball of \mathbb{R}^{4^n-1} , such that $\varphi(x) = 0$. Then for a given $x_0, x_1 \in P_x$, and any Finsler metric $F(x, y)$, there exist a constant minimum $\mathfrak{m} > 0$ and a constant maximum $\mathfrak{M} > 1$ such that the following relation is true:*

$$\mathfrak{m} |\varphi(x_1) - \varphi(x_0)| \leq L_F(x_0, x_1) \leq \mathfrak{M} |\varphi(x_1) - \varphi(x_0)| . \quad (3)$$

Here $|\varphi(x_1) - \varphi(x_0)|$ denotes the Euclidean length of the $4^n - 1$ dimensional vector in the tangent space.

Proof: We first note that a compact set P_x for which $\varphi(x) = 0$ always exists. This is true because, given a local coordinate system $\varphi : Q \rightarrow \mathbb{R}^{4^n-1}$ and $x \in Q$ for which $\varphi(x) = 0$, we can choose P_x to be a closure of the preimage of $\varphi^{-1}(B^{4^n-1}(r))$ for some $r > 0$. By $B^{4^n-1}(r) = \{v \in \mathbb{R}^{4^n-1} : |v| = \sqrt{\sum_i v_i^2} < r\}$ we denote the ball of radius r in the tangent space whose closure is a subset of $\varphi(Q)$.

Next we note that, for tangent vector $y = \sum_i y_i \frac{\partial}{\partial x_i} \equiv \frac{dx}{dt} \in T_x \mathcal{U}$, the ratio between Minkowski norm $F(x, y)$ and x -dependent Euclidean norm $|y| := \sqrt{\sum_i y_i^2}$ for the basis $\{\frac{\partial}{\partial x_i}\}$ is well defined for $y \neq 0$. Since both norms are positive continuous functions over the compact sets their quotient is also a positive continuous function. Therefore the quotient's minimum \mathfrak{m} and maximum \mathfrak{M} exist and are both positive: $0 < \mathfrak{m} \leq \frac{F(y)}{|y|} \leq \mathfrak{M}$. In other words, for all $y \in T_x \mathcal{U}$ and all $x \in P_x$:

$$\mathfrak{m}|y| \leq F(x, y) \leq \mathfrak{M}|y|. \quad (4)$$

Now we can prove the right hand side (RHS) of inequality (4) by choosing the path $\sigma \in P$ that maps under φ to a line segment. In that case we can write:

$$L_F(x_0, x_1) = \int_{t_0}^{t_1} F(\sigma') dt \leq \mathfrak{M} \int_{t_0}^{t_1} |\sigma'| dt = \mathfrak{M} |\varphi(x_1) - \varphi(x_0)|, \quad (5)$$

where $\sigma' = \frac{d\sigma}{dt}$ denotes the velocity field of a path σ .

To prove the left hand side of inequality (4) we first show that σ must be contained in P_x . The proof is by contradiction as follows.

Choose $r_0 < \frac{\mathfrak{m}}{\mathfrak{m}+3\mathfrak{M}}r$ and $\epsilon_0 = \mathfrak{M}r_0$ and $P_0 = \varphi^{-1}[B^n(r_0)] \subset P_x$. Let $\sigma : [t_0, t_1] \rightarrow \mathcal{U}$ be a piecewise C^∞ curve such that $\sigma(t_0) = x_0$ and $\sigma(t_1) = x_1$ for $x_0, x_1 \in P_0$. If $L_F(\sigma) \leq d_F(x_0, x_1) + \epsilon_0$ then the curve σ is certainly contained in P_x , and since by equation (5) $d_F(x_0, x_1) \leq 2\mathfrak{M}r_0$ we have by assumption that $L_F(\sigma) \leq 3\mathfrak{M}r_0$. Now if we suppose that σ is not contained in P_x , and let $t_0 \leq t^* \leq t_1$ be the first instance where σ reaches the boundary ∂P_x , at the point $q \equiv \sigma(t^*)$, so that $|\varphi(q)| = r$, then:

$$L_F(\sigma) \geq L_F(\sigma_{[t_0, t^*]}) = \int_{t_0}^{t^*} F(\sigma') dt \geq \mathfrak{m} \int_{t_0}^{t^*} |\sigma'| dt \geq \mathfrak{m} |\varphi(q) - \varphi(x_0)| \geq \mathfrak{m}(r - r_0). \quad (6)$$

But the length of this curve would in fact be longer than the maximum possible length of $3\mathfrak{M}r_0 < \mathfrak{m}(r - r_0)$, since by assumption we are assured that for $\mathfrak{m} > 0$ and $\mathfrak{M} > 1$ it is true that $r_0 < \frac{\mathfrak{m}}{\mathfrak{m}+3\mathfrak{M}}r$. Therefore σ must be contained in P_x .

The proof of the left hand side of inequality (3) follows by the same arguments as were used to prove (6)□

Given the distortion lemma for the length of the path for any two points that belong to the compact set, we can easily derive a similar result that is valid for the shortest distances.

Corollary 1. *For a Finsler manifold (\mathcal{U}, F) , and any point $x \in \mathcal{U}$, let $\varphi : P_x \rightarrow \mathbb{R}^{4^n-1}$ be the local coordinate system diffeomorphism of a compact set P_x onto an open ball of \mathbb{R}^{4^n-1} , such that $\varphi(x) = 0$. Then, for a given $x_0, x_1 \in P_x$ and any Finsler metric $F(x, y)$ there exist a constant minimum $\mathfrak{m} > 0$ and a constant maximum $\mathfrak{M} > 1$ such that the following relation is true:*

$$\mathfrak{m}|\varphi(x_1) - \varphi(x_0)| \leq d_F(x_0, x_1) \leq \mathfrak{M}|\varphi(x_1) - \varphi(x_0)|. \quad (7)$$

Proof: We only need to verify the left hand side of inequality (3) is still true for minimal length curves. By definition of metric distance, for $0 \leq \epsilon \leq \epsilon_0$, two points $x_0, x_1 \in P_0$ can be joined by a piecewise C^∞ curve $\sigma : [t_0, t_1] \rightarrow \mathcal{U}$ with integral length:

$$L_F(\sigma) \leq d_F(x_0, x_1) + \epsilon . \quad (8)$$

By previous arguments, σ must lie in P_x , end by similar calculations we find that:

$$\mathfrak{m}|\varphi(x_1) - \varphi(x_0)| \leq L_F(\sigma) \leq d_F(x_0, x_1) + \epsilon , \quad (9)$$

Letting $\epsilon \rightarrow 0$ proves the desired result. \square

Lemma (1) and Corollary (1) allow us to bound the lengths on the manifold to the Euclidian lengths on the tangent space. For Euclidean coordinates in our tangent space we will have the coefficients in the decomposition of the gate Hamiltonian matrix in terms of the generalized Pauli matrices, n times tensored two dimensional matrices from the set $\{I, X, Y, Z\}$.

For example, in the context of simulation, points $x_0, x_1 \in P_x$ and an open set P_x are chosen such that they correspond to the end points in the simulation by a single gate. Moreover, we can construct a *local coordinate system* on the Lie group $SU(2^n)$ which is a Lie algebra $\mathfrak{su}(2^n)$. For the *origin* $x_s \in \mathcal{U}$, define a pull back map $\varphi^{-1} : \mathbb{R}^{4^n-1} \rightarrow \mathcal{U}$, so that $x_{s+1} \equiv \exp^{-iy_{s+1} \cdot \sigma} x_s = \exp^{-i\varphi(x_{s+1}) \cdot \sigma} x_s$, where σ denotes the coordinate basis, i.e. $(4^n - 1)$ -component vector whose entries are the generalized Pauli matrices.

For the particularly chosen metric, as in [4], $F_p(x_0, y) \equiv F_p(y) = \sqrt{\sum_{i=1}^k y_i^2 + p^2 \sum_{j \neq i} y_j^2}$, where $k = \frac{9(n^2-n)}{2} + 3n$, which introduces a penalty p for the subset of Hamiltonian coordinates in the tangent space, so we have:

$$|y| \leq F_p(y) \leq p|y|.$$

Since this relation is true on any compact set, by Corollary (1) we have:

$$|\varphi(x_{s+1}) - \varphi(x_s)| \leq d_{F_p}(x_s, x_{s+1}) \leq p|\varphi(x_{s+1}) - \varphi(x_s)|. \quad (10)$$

It is important to note that in our analysis constants \mathfrak{m} and \mathfrak{M} do not depend on the compact set within which each gate is applied, and they are basically the property of the metric. This property might not be true in general for some other Finsler structures, but for our purposes here this assumption is very plausible.

4 Equivalence Result

For the sake of consistency and easier understanding, we follow the notation from [3] and denote with $m_{\mathcal{G}}$ the minimum number of gates, for a given set \mathcal{G} , needed to implement an arbitrary unitary $U \in \mathcal{U}$. Moreover, in this section we assume that the geodesic is simulated by sequential application of the gates from the set \mathcal{G} , and that by using a single gate from the set \mathcal{G} we can simulate exactly any other point in the ball $B_{x_0}^+(\epsilon)$, i.e. which is the ϵ -neighborhood around the initial condition at the point x_0 . This is an unrealistic scenario, since the set of gates would need to be infinite and non-local. Hence, we relax it in the next section. However, for the purpose of exact simulation of the geodesic it is an important tool. Clearly consideration of the set of gates defined in this way, as we shall see, is the best we can possibly hope for, and thus the bounds achieved by the set \mathcal{G} are optimal, i.e. they determine the bounds achievable by any other set of gates that is less powerful.

For any gate used in a simulation we assign a gate index, so that eventually the index set is $s = \{0, 1, 2, \dots, m_{\mathcal{G}} - 1\}$ for every gate in the simulation. Moreover, by $\sigma(t) : [0, m_{\mathcal{G}}] \rightarrow \mathcal{U}$ we denote a minimal geodesic between $\sigma(0) = I$, $\sigma(m_{\mathcal{G}}) = U$.

Note that, since \mathcal{U} is a compact manifold with Finsler structure, all forward and backward Cauchy sequences with respect to d must converge on \mathcal{U} . More precisely, compact Finsler spaces are automatically both forward complete and backward complete. This fact holds regardless of whether the Finsler structure is absolutely homogeneous or only positively homogeneous. Therefore, any two points on the manifold can be connected by a minimizing geodesic, as that property itself is a sufficient condition for the Hopf-Rinow theorem [6].

Theorem 1. *Let $d_F(I, U)$ denote a length of a geodesic between I and $U \in SU(2^n)$. For any simulation index set $s = \{0, 1, 2, \dots, m_{\mathcal{G}} - 1\}$ let $P_{x_s} \in \mathcal{U}$ be an open set on the manifold that contains a segment of minimizing geodesic $\sigma_s(t) : [s, s+1] \rightarrow \mathcal{U}$, that is simulated exactly by a single gate. Moreover, let P_{x_s} be mapped by φ diffeomorphically onto an open ball in \mathbb{R}^{4^n-1} , so that $\rho_s = |\varphi(x_{s+1}) - \varphi(x_s)|$ is the Euclidean length of the image of the geodesic segment σ_s . If we denote $\rho_{\sup} = \sup_s \rho_s$ and $\rho_{\inf} = \inf_s \rho_s$, then the following relation holds:*

$$\frac{d_F(I, U)}{\rho_{\sup} \mathfrak{M}} \leq m_{\mathcal{G}} \leq \frac{d_F(I, U)}{\rho_{\inf} \mathfrak{m}} . \quad (11)$$

Proof: For any segment gate index from set s , by the Corollary (1) we see that:

$$\mathfrak{m} \rho_s \leq d_F(x_s, x_{s+1}) \leq \mathfrak{M} \rho_s ,$$

Summing over all segments of minimizing geodesic $\sum_{s=0}^{m_{\mathcal{G}}-1} d_F(x_s, x_{s+1}) = d_F(I, U)$, and taking into account that $\sum_{s=0}^{m_{\mathcal{G}}-1} \beta_s \leq m_{\mathcal{G}} \beta$ and $m_{\mathcal{G}} \frac{\rho_s^2}{\beta} \leq \sum_{s=0}^{m_{\mathcal{G}}-1} \frac{\rho_s^2}{\beta_s}$, it is easy to see that:

$$m_{\mathcal{G}} \mathfrak{m} \rho_{\inf} \leq d_F(I, U) = \sum_{s=0}^{m_{\mathcal{G}}-1} d_F(x_s, x_{s+1}) \leq m_{\mathcal{G}} \mathfrak{M} \rho_{\sup} , \quad (12)$$

which gives the desired result by rearranging the variables. \square

Note that the above theorem is derived in terms of bounds of the Euclidean distances in the tangent space. One may take a different path though, as for example Nielsen in [3], by deriving the result for the lower bound in terms of the lengths of the geodesic segments simulated by the single gate: $d_F(x_s, x_{s+1}) \leq \beta_{\sup}$. From the following theorem, one can reproduce the result derived by Nielsen as a special case when $\beta_{\sup} = 1$.

Theorem 2. *Let $d_F(I, U)$ denote a length of a geodesic between I and $U \in SU(2^n)$. For any simulation index set $s = \{0, 1, 2, \dots, m_{\mathcal{G}} - 1\}$ let $P_{x_s} \in \mathcal{U}$ be an open set on the manifold that contains a segment of minimizing geodesic $\sigma_s(t) : [s, s+1] \rightarrow \mathcal{U}$, that is simulated exactly by a single gate. Moreover, let P_{x_s} be mapped by φ diffeomorphically onto an open ball in \mathbb{R}^{4^n-1} , so that $\rho_s = |\varphi(x_{s+1}) - \varphi(x_s)|$ is an image of the bounded length geodesic segment $d_F(x_s, x_{s+1}) \leq \beta_s$. If we denote $\beta_{\sup} = \sup_s \beta_s$ and $\beta_{\inf} = \inf_s \beta_s$, then the following relation holds:*

$$\frac{d_F(I, U)}{\beta_{\sup}} \leq m_{\mathcal{G}} \leq \frac{\mathfrak{M}}{\mathfrak{m}} \frac{d_F(I, U)}{\beta_{\inf}} . \quad (13)$$

Proof: Following along the lines of Theorem (1):

$$\mathfrak{m} \frac{\beta_s}{\mathfrak{M}} = \mathfrak{m} \rho_s \leq d_F(x_s, x_{s+1}) \leq \beta_s \equiv \mathfrak{M} \rho_s ,$$

Summing over all segments of minimizing geodesic $\sum_{s=0}^{m_g-1} d_F(x_s, x_{s+1}) = d_F(I, U)$, and taking into account that $m_{\mathcal{G}}\beta_{\inf} \leq \sum_{s=0}^{m_g-1} \beta_s \leq m_{\mathcal{G}}\beta_{\sup}$:

$$\frac{m}{M} m_{\mathcal{G}}\beta_{\inf} \leq d_F(I, U) \leq m_{\mathcal{G}}\beta_{\sup} , \quad (14)$$

which gives the stated result. \square

Equations (11) and (14) establish the tightest possible equivalence between the minimal number of gates in the circuit and geodesic length as a function of the simulation parameters. Again, the simulation parameters may be defined in terms of distances traversed with the single gate on the manifold or in terms of the Euclidean distances between the initial and final coefficients in the generalized Pauli expansion of the gate Hamiltonian. Even though the above results give no indication as to how to implement the simulation, they do provide us the best bounds we currently have and give us an estimate to the quality of the simulation provided that one knows the simulation parameters. However, the above results can be applied to the arbitrary paths on the manifold including those that are generated adiabatically. In particular, it would be very interesting to compare the results for bounds of circuit size obtained by geometric techniques with the equivalence results obtained in [1].

5 Approximate simulation

In this section we reformulate the bounds for the standard circuit simulation procedure where the set of gates used consists solely of the single and two qubit gates, which are applied sequentially. Since the exact simulation of arbitrary unitary gate by single and two qubit gates demands an exponential number of gates, almost all unitaries simulated by the polynomial number of gates will be simulated approximately.

In particular, we consider two paths. Let the first be $d_{F_p}(I, \tilde{U})$, denoting the length of the geodesic simulated exactly with the set of gates from \mathcal{G} with respect to the Finsler metric F_p , and let the second one $L_{F_p}(I, \tilde{U})$ be the minimum length path for the exact simulation of \tilde{U} by the set of gates from \mathcal{G}_2 . Here we denote by \mathcal{G}_2 the set of unitary gates whose time independent Hamiltonians have Hamming weight not greater than two.

Note that the length $L_{F_p}(I, \tilde{U})$ has nothing to do with $d_{F_p}(I, \tilde{U})$, as $L_{F_p}(I, \tilde{U})$ is completely determined by the simulation, and almost everywhere does not simulate the geodesic $d_{F_p}(I, \tilde{U})$.

Corollary 2. *Let \tilde{U} be the approximation of the unitary operation U that is simulated by the one and two qubit gates. Then the lower bound on the minimum circuit size $\tilde{m}_{\mathcal{G}_2}$ is at most $O(\frac{n^4}{p} d_{F_p}^2(I, U) L_{F_p}(I, \tilde{U}))$, and the upper bound on $m_{\mathcal{G}_2}$ is at least $\Omega(n^4 d_{F_p}^3(I, U))$.*

Proof: The three step standard simulation of arbitrary $U = e^{-iH(t)t}$ is elaborated in detail by Nielsen, Dowling, Gu and Doherty in [4]. The procedure can be sketched as follows:

- (1) the time variable Hamiltonian $H(t)$ is substituted by projected the Hamiltonian $H_p(t)$ that is formed by deleting all σ_i for $i > k$, i.e. all three- and more-body terms in the Pauli expansion of $H(t) = \sum_{i=1}^k y_i \sigma_i + \sum_{j \neq i} y_i \sigma_i$, where $k = \frac{9(n^2-n)}{2} + 3n$;
- (2) the evolution due to $H_p(t)$ is broken up into many small intervals, each of length Δ , over which the time-dependent Hamiltonian $H_p(t)$ is accurately simulated by a constant mean Hamiltonian \bar{H}_p^Δ ;

- (3) the mean Hamiltonian \bar{H}_p^Δ that has k terms in the Pauli expansion with coefficients $|y_i| \leq 1$ is simulated with a standard simulation technique [14] using one and two qubit gates.

The reader is encouraged to see [4] for full detail of the approximation result.

For the above procedure, since $SU(2^n)$ is compact and simply connected, there exists a path $L_{F_p}(I, \tilde{U})$ that is exactly synthesized with the gates in the simulation. By exactly simulated we mean that the end points of each gate in the simulation lie precisely on the path of length $L_{F_p}(I, \tilde{U})$. Clearly, the length of $L_{F_p}(I, \tilde{U}) \geq d_{F_p}(I, U)$.

Now we bound length of the path segments, $L_{F_p}(\tilde{x}_s, \tilde{x}_{s+1})$, for each of $\tilde{m}_{\mathcal{G}_2}$ gates in the simulation. Since there exists a compact set P_{x_s} , such that end points $\tilde{x}_s, \tilde{x}_{s+1} \in P_{x_s}$, that maps diffeomorphically to the local coordinate system, we can use Lemma (1) and its corollaries. Corroborating the arguments used to derive equation (10), over the compact set P_{x_s} , the Finsler structure, i.e. the Minkowski norm for the Pauli expansion of $H(t)$, is $F_p(x_s, y_s) = \sqrt{\sum_{i=1}^k y_i^2 + p^2 \sum_{j \neq i} y_j^2}$. Its minimum and maximum distortion over the compact set P_{x_s} are: $|y_s| \leq F_p(y_s) \leq p|y_s|$. Therefore, by the Lemma (1)

$$|\varphi(\tilde{x}_{s+1}) - \varphi(\tilde{x}_s)| \leq L_{F_p}(\tilde{x}_s, \tilde{x}_{s+1}) \leq p|\varphi(\tilde{x}_{s+1}) - \varphi(\tilde{x}_s)|.$$

The same is true for any other segment in the simulation, and hence:

$$m_{\mathcal{G}_2} \rho_{\inf}^\Delta \leq L_{F_p}(I, \tilde{U}) = \sum_{s=0}^{m_g-1} L_{F_p}(\tilde{x}_s, \tilde{x}_{s+1}) \leq m_{\mathcal{G}_2} p \rho_{\sup}^\Delta \quad (15)$$

where $\rho_{\inf}^\Delta = \inf_s |\varphi(\tilde{x}_{s+1}) - \varphi(\tilde{x}_s)|$, and $\rho_{\sup}^\Delta = \sup_s |\varphi(\tilde{x}_{s+1}) - \varphi(\tilde{x}_s)|$. Note that we can always choose the s -th gate local coordinate system so that $\varphi(\tilde{x}_s) = 0$.

Finally, in the three-step simulation summarized above, gates at the third stage simulate the time invariant Hamiltonian \bar{H}_p^Δ for the segment Δ , with coordinates $|y_i| \leq 1$. More precisely, the s -th gate simulates the neighborhood around x_s : $x_{s+1} \equiv e^{-i\varphi(x_{s+1}) \cdot \sigma} x_s = e^{-iy_s \sigma_s \Delta^2} x_s$. Here $\sigma \in \mathcal{G}$ denotes stabilizer basis on n qubits, $\sigma_s \in \mathcal{G}_2$, and Δ^2 is the simulation time for every gate. If we choose $\Delta = \Theta((n^2 d_{F_p}(I, U))^{-1})$, as in [4], then for $|y_s| \leq 1$ we see that $\rho_s = |y_s \Delta^2| = \Theta((n^4 d_{F_p}^2(I, U))^{-1})$. Finally, using equation (15) we establish that the simulation with $\tilde{m}_{\mathcal{G}_2}$ gates has the upper bound $\Theta(n^4 d_{F_p}^2(I, U) L_{F_p}(I, \tilde{U})) \geq \Omega(n^4 d_{F_p}^3(I, U))$. By similar arguments, for the lower bound, we get $O(\frac{n^4}{p} d_{F_p}^2(I, U) L_{F_p}(I, \tilde{U}))$. \square

6 Conclusion

The Distortion Lemma and its corollary provide a general tool for relating distances on the manifold with distances on the tangent space. In this paper we have derived a generalized linear bounds for the exact simulation of any path on the manifold, in terms of the minimum circuit size and the simulation parameters.

The equivalence between the path on the manifold and circuit size still persists in the case of approximate simulation, provided that the simulation parameters have the appropriate scaling. However, one can not expect better than n^2 times improvement in the minimum circuit size upper bound over the result for standard circuit simulation derived by Nielsen, Dowling, Gu and Doherty [4].

Moreover, if one defines a metric on the manifold that penalizes the hard-to-simulate directions on the tangent space with high cost, that cost is, in effect, translated to the increased ratio between upper and lower bound in minimum circuit size.

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